



STIC Search Report

Biotech-Chem Library

STIC Database Tracking Number: 199228

TO: Shailendra Kumar
Location: REM-5C03/5C18
Art Unit: 1621
Friday, August 25, 2006
Case Serial Number: 10/522077

From: Usha Shrestha
Location: Biotech-Chem Library
REM-1A64
Phone: (571)272-3519

Usha.shrestha@uspto.gov

Search Notes

Examiner Kumar,

See attached results.

If you have any questions about this search feel free to contact me at any time.

Thank you for using STIC search services!

Usha Shrestha
Technical Information Specialist
STIC Biotech/Chem Library
(571)272-3519

8-999

199228

Access DB# _____

SEARCH REQUEST FORM

Scientific and Technical Information Center

Requester's Full Name: S. Kumar Examiner #: 695-914 Date: 8/22/06
 Art Unit: 1621 Phone Number: 301-20640 Serial Number: 101522077
 Mail Box and Bldg/Room Location: REM 5C03 Results Format Preferred (circle): PAPER DISK E-MAIL
5C18

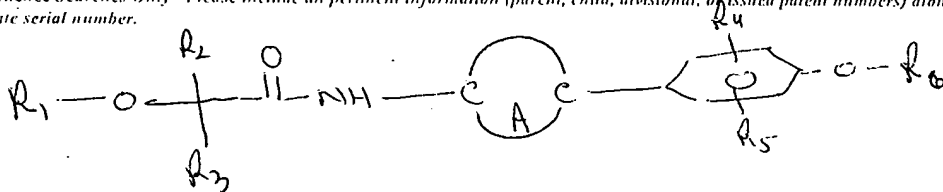
If more than one search is submitted, please prioritize searches in order of need.

Please provide a detailed statement of the search topic, and describe as specifically as possible the subject matter to be searched. Include the elected species or structures, keywords, synonyms, acronyms, and registry numbers, and combine with the concept or utility of the invention. Define any terms that may have a special meaning. Give examples or relevant citations, authors, etc, if known. Please attach a copy of the cover sheet, pertinent claims, and abstract.

Title of Invention: N-bisaryl- and n-aryl cycloalkylidenyl - alpha-hydroxy-
 Inventors (please provide full names): Clemens Lamberth et al.

Earliest Priority Filing Date: 7/24/02

For Sequence Searches Only Please include all pertinent information (parent, child, divisional, or issued patent numbers) along with the appropriate serial number.



R_1, R_2, R_4, R_5 are H, alkyl, ~~alkenyl~~, alkynyl, etc.
 R_6 is propargyl
 A is 1,2-cyclohexylidene or 1,2-cyclopropylidene

Also see claims 10 + 11 + 12 for process

STAFF USE ONLY

Type of Search

Vendors and cost where applicable

Searcher: UAC

NA Sequence (#) _____

STN 348-15

Searcher Phone #: _____

AA Sequence (#) _____

Dialog _____

Searcher Location: _____

Structure (#) 4

Questel/Orbit _____

Date Searcher Picked Up 8/24/06

Bibliographic _____

Dr. Link _____

Date Completed: 8/25/06

Litigation _____

Lexis/Nexis _____

Searcher Prep. Review Time: 60

Fulltext _____

Sequence Systems _____

Clerical Prep. Time: 30

Patent Family _____

WWW/Internet _____

Online Time 120

Other _____

Other (specify) _____

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FILE 'REGISTRY' ENTERED AT 10:02:01 ON 25 AUG 2006

=> d his

FILE 'HCAPLUS' ENTERED AT 08:04:31 ON 25 AUG 2006

L1 1 S US20050245607/PN
SEL RN

FILE 'REGISTRY' ENTERED AT 08:04:45 ON 25 AUG 2006

L2 75 S E1-E75

FILE 'REGISTRY' ENTERED AT 08:18:47 ON 25 AUG 2006

L3 STR
L4 0 S L3
L5 STR L3
L6 1 S L5
L7 STR
L8 STR
L9 3 S L7 AND L8
L10 STR L7
L11 STR L7
L12 3 S L10 AND L11 AND L8
L13 2 S L12 AND L2
L14 35 S L10 AND L11 AND L8 FUL
L15 24 S L14 AND L2
SAV L14 KUM077/A

FILE 'HCAPLUS' ENTERED AT 09:56:08 ON 25 AUG 2006

L16 3 S L14

FILE 'CASREACT' ENTERED AT 09:58:48 ON 25 AUG 2006

L17 STR L3
L18 0 S L17 SAM
L19 0 S L17 FUL

=> d que l16

L8 STR

O~C~C~O
1 2 3 4

NODE ATTRIBUTES:

DEFAULT MLEVEL IS ATOM

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 4

STEREO ATTRIBUTES: NONE

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CH2-C \equiv C
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L16 3 SEA FILE=HCAPLUS ABB=ON L14

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FILE 'HCAPLUS' ENTERED AT 10:02:25 ON 25 AUG 2006

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L16 ANSWER 1 OF 3 HCAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2004:101121 HCAPLUS
DOCUMENT NUMBER: 140:163585
TITLE: Preparation of N-bisaryl- and
N-aryl-cycloalkylidenyl-alpha-hydroxy- and
alpha-alkoxy acid amides as fungicides
INVENTOR(S): Lamberth, Clemens; Zeller, Martin; Goegh,
Tibor
PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.
SOURCE: PCT Int. Appl., 70 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004011417	A1	20040205	WO 2003-EP8057	2003 0723

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA,
CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI,
GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG,
KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK,
MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU,

SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA,
 UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM,
 AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ,
 DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL,
 PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN,
 GQ, GW, ML, MR, NE, SN, TD, TG

CA 2493975 AA 20040205 CA 2003-2493975

2003
 0723

AU 2003251456 A1 20040216 AU 2003-251456

2003
 0723

EP 1534664 A1 20050601 EP 2003-771074

2003
 0723

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE,
 MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ,
 EE, HU, SK

BR 2003012872 A 20050614 BR 2003-12872

2003
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CN 1671651 A 20050921 CN 2003-817601

2003
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JP 2005533854 T2 20051110 JP 2004-523767

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US 2005245607 A1 20051103 US 2005-522077

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PRIORITY APPLN. INFO.:

GB 2002-17211

A

2002
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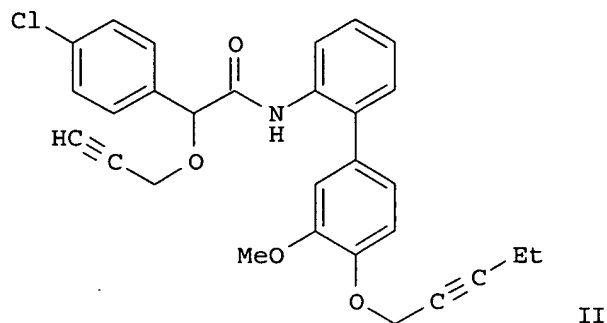
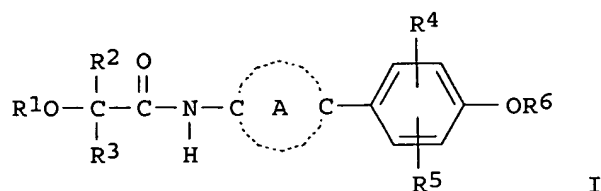
WO 2003-EP8057

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OTHER SOURCE(S):
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MARPAT 140:163585



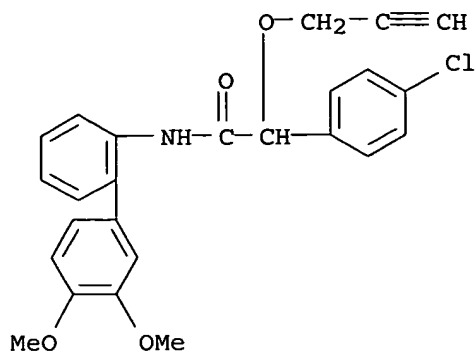
AB The title compds. [I; R1 = H, alkyl; alkenyl, alkynyl, haloalkyl; R2 = H, (un)substituted alkyl, alkenyl, alkynyl; R3 = (un)substituted (hetero)aryl; A = (un)substituted (un)saturated cycloalkylidene, phenylidene or (un)saturated heterocyclylidene bridge; R4, R5 = H, an organic radical; R6 = H, trialkylsilyl, dialkylphenylsilyl, alkyldiphenylsilyl, triphenylsilyl, (un)substituted alkyl, alkenyl or alkynyl], were prepared E.g., a multi-step synthesis of II (starting from 4-bromoguaiacol), was given. The compds. I possess plant-protecting properties and are suitable for protecting plants against infestation by phytopathogenic microorganism, especially fungi. E.g., some of them inhibited fungal infestation to at least 80% in three different tests.

IT 602280-76-2P 602280-77-3P 655222-80-3P
 655222-81-4P 655222-82-5P 655223-00-0P
 655223-03-3P 655223-06-6P 655223-12-4P
 655223-13-5P 655223-14-6P 655223-15-7P
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 655223-19-1P 655223-20-4P 655223-21-5P
 655223-22-6P 655223-23-7P 655223-24-8P
 655223-25-9P 655223-26-0P

(preparation of N-aryl- α -hydroxy(or alkoxy) phenylacetamides as fungicides)

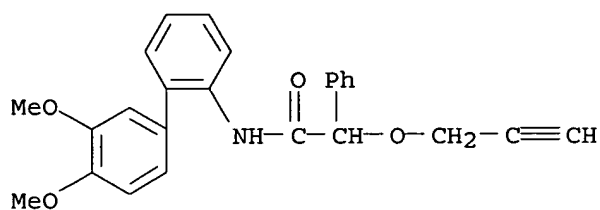
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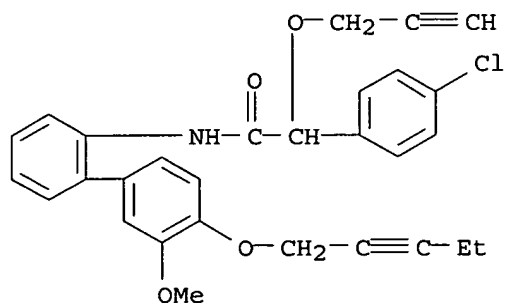
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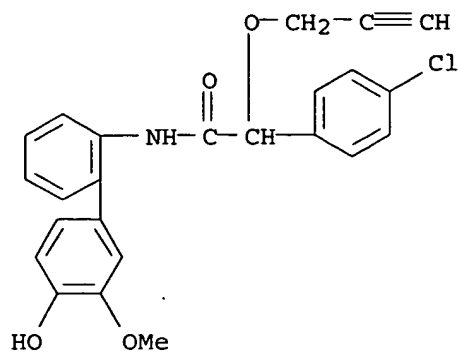
RN 655222-80-3 HCAPLUS

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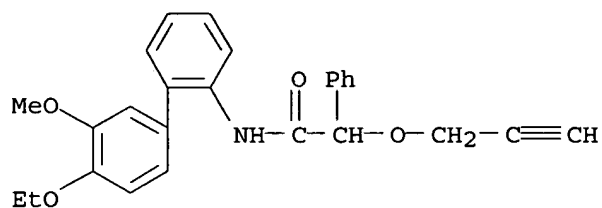


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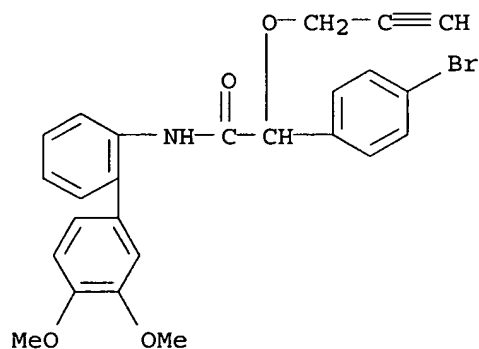
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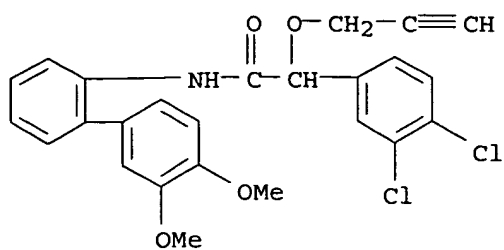
CN Benzeneacetamide, N-(4'-ethoxy-3'-methoxy[1,1'-biphenyl]-2-yl)-
α-(2-propynyloxy) - (9CI) (CA INDEX NAME)

RN 655223-00-0 HCAPLUS

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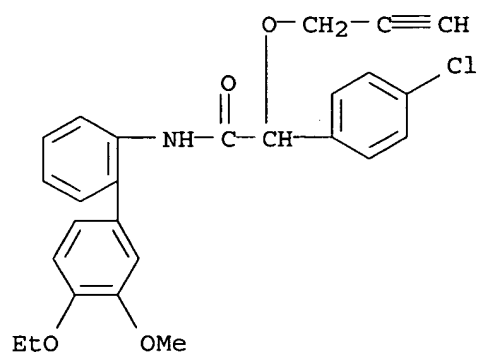
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CN Benzeneacetamide, 3,4-dichloro-N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-
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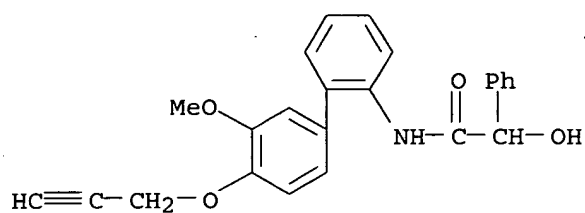
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CN Benzeneacetamide, 4-chloro-N-(4'-ethoxy-3'-methoxy[1,1'-biphenyl]-2-yl)-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



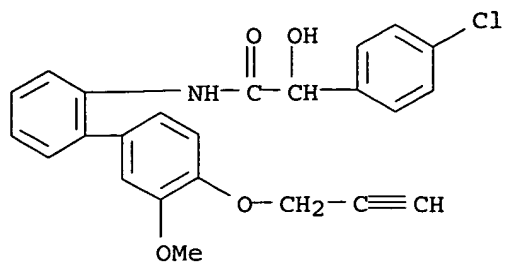
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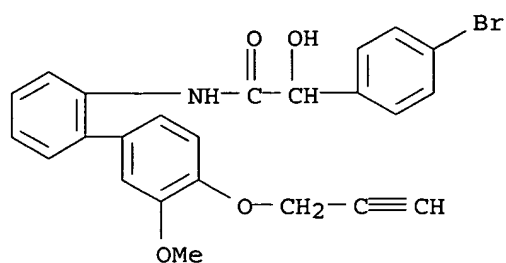
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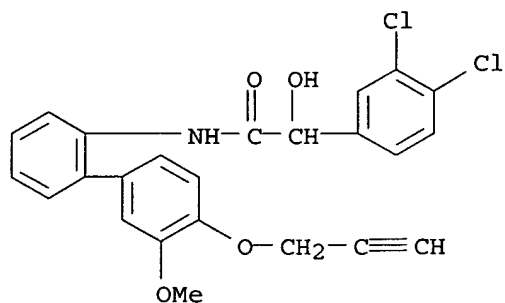
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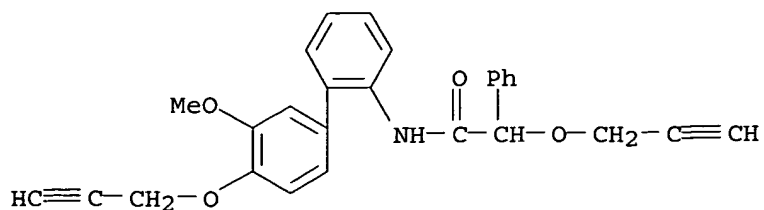
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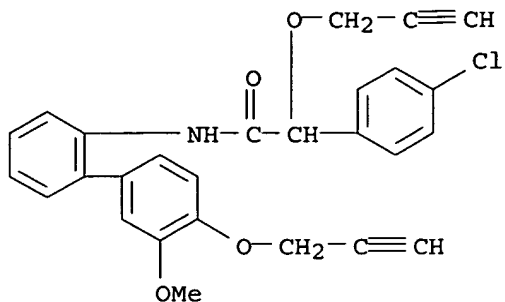


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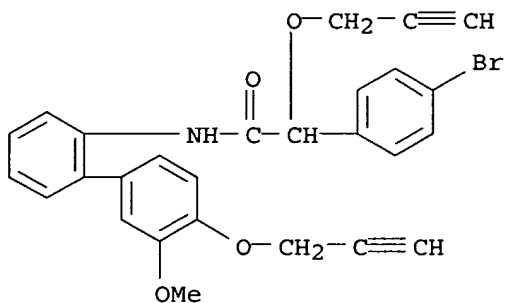
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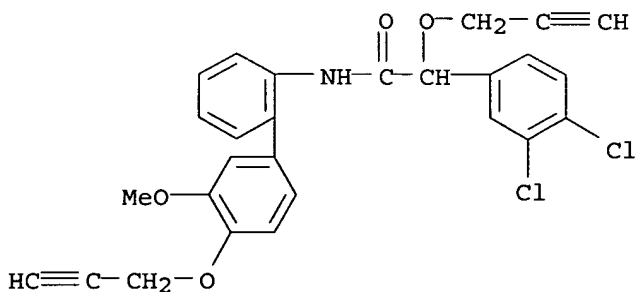
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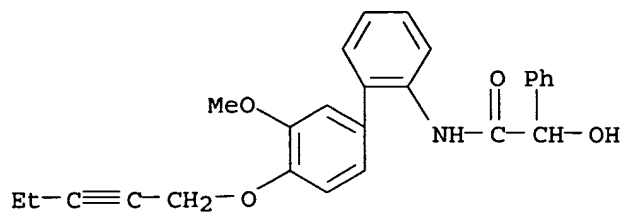
CN Benzeneacetamide, 4-bromo-N-[3'-methoxy-4'-(2-propynyloxy) [1,1'-biphenyl]-2-yl]- α -(2-propynyloxy) - (9CI) (CA INDEX NAME)

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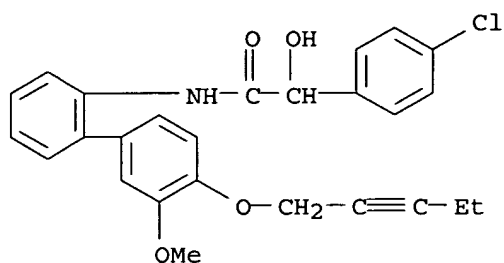
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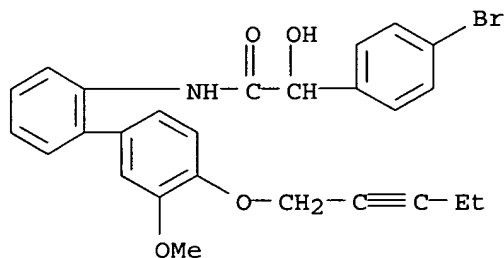
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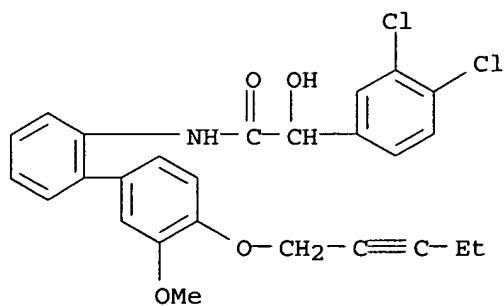
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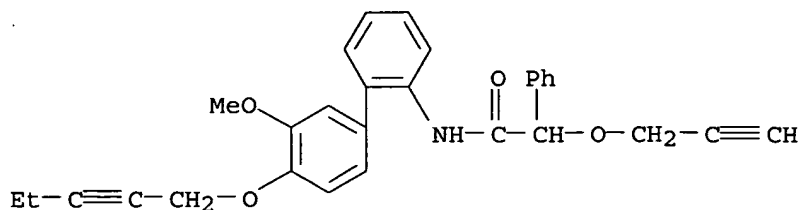


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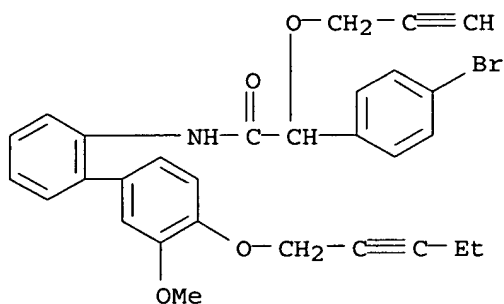
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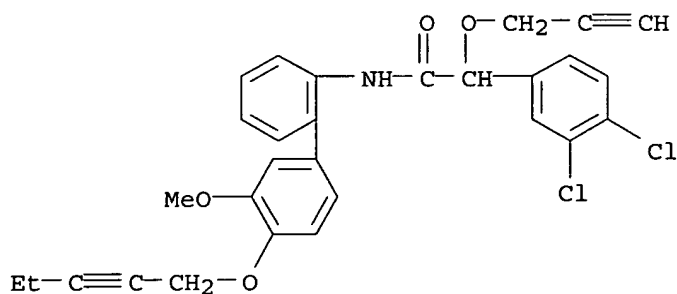
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RN 655223-25-9 HCAPLUS

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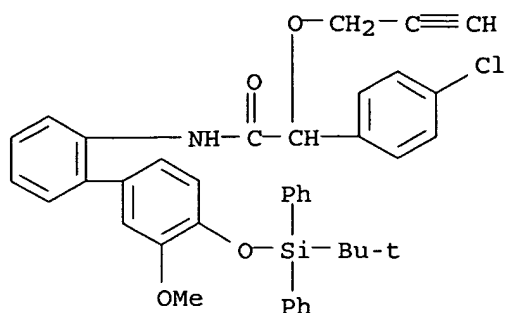
CN Benzeneacetamide, 3,4-dichloro-N-[3'-methoxy-4'-(2-pentynyloxy) [1,1'-biphenyl]-2-yl]- α -(2-propynyloxy) - (9CI) (CA INDEX NAME)

IT 655223-08-8P

(preparation of N-aryl- α -hydroxy(or alkoxy) phenylacetamides as fungicides)

RN 655223-08-8 HCAPLUS

CN Benzeneacetamide, 4-chloro-N-[4'-[[[(1,1-dimethylethyl)diphenylsilyl]oxy]-3'-methoxy[1,1'-biphenyl]-2-yl]- α -(2-propynyloxy) - (9CI) (CA INDEX NAME)



IC ICM C07C235-38
ICS C07C235-36; C07C215-74; C07C217-80; C07C215-64; C07C217-74

CC 25-19 (Benzene, Its Derivatives, and Condensed Benzenoid Compounds)

Section cross-reference(s): 5

IT 602280-76-2P 602280-77-3P 602280-96-6P
602281-04-9P 655222-80-3P 655222-81-4P
655222-82-5P 655222-83-6P 655222-84-7P 655222-85-8P
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655223-35-1P 655223-36-2P 655223-37-3P
(preparation of N-aryl- α -hydroxy(or alkoxy) phenylacetamides as fungicides)
IT 431879-38-8P 482309-54-6P 482309-55-7P 482309-56-8P
482309-57-9P 482309-58-0P 655223-08-8P
(preparation of N-aryl- α -hydroxy(or alkoxy) phenylacetamides as fungicides)

REFERENCE COUNT: 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L16 ANSWER 2 OF 3 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:757666 HCAPLUS

DOCUMENT NUMBER: 139:261052

TITLE: Preparation of N-biphenylacetamides as agrochemical fungicides

INVENTOR(S): Sakaguchi, Hiroshi

PATENT ASSIGNEE(S): Sumitomo Chemical Company, Limited, Japan

SOURCE: PCT Int. Appl., 89 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: Japanese

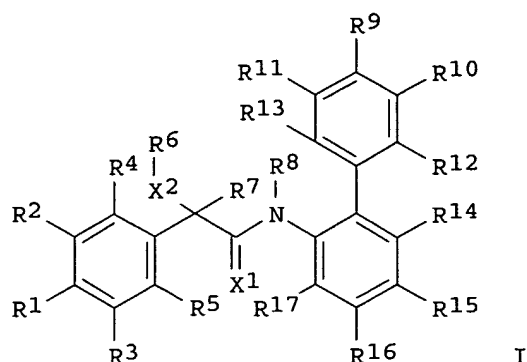
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PATENT INFORMATION:

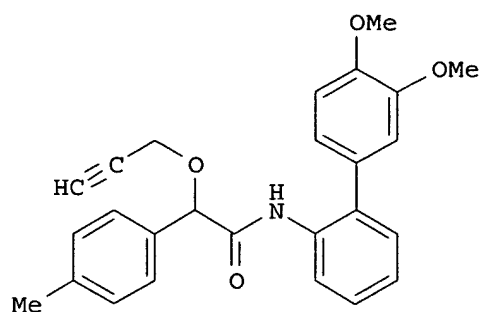
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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USHA SHRESTHA EIC 1600 REM 1A64

WO 2003078382	A1	20030925	WO 2003-JP2015	2003 0225
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, KE, KG, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MX, MY, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2003211652	A1	20030929	AU 2003-211652	2003 0225
EP 1486483	A1	20041215	EP 2003-707030	2003 0225
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
JP 2003335744	A2	20031128	JP 2003-67817	2003 0313
US 2005192359	A1	20050901	US 2004-507586	2004 0913
RITY APPLN. INFO.:			JP 2002-71627	A 2002 0315
			WO 2003-JP2015	W 2003 0225
R SOURCE(S) :	MARPAT 139:261052			



I



II

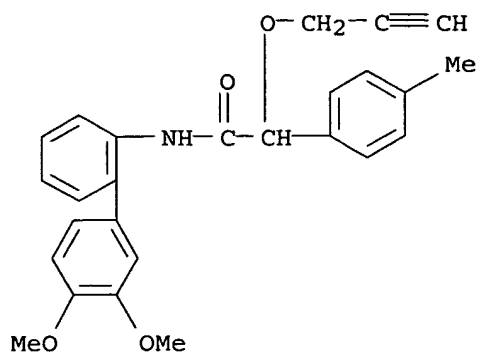
AB The title N-biphenylacetamides with general formula of I [R1-R5 = independently H, halo, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)alkoxy, (halo)alkenyloxy, (halo)alkynyloxy, (halo)alkylthio, cycloalkyl(oxy), trialkylsilyl, or CN; R6 = (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)alkylcarbonyl, (halo)alkenylcarbonyl, or (halo)alkynylcarbonyl; R7 = H or alkyl; R8 = H or alkyl; R9 and R10 = independently halo, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, cyanoalkyl, (halo)alkoxy, (halo)alkenyloxy, (halo)alkynyloxy, cyanoalkoxy, (halo)alkylthio, alkenylthio, alkynylthio, cycloalkyl(oxy), cycloalkylalkoxy, alkoxyalkyl, trialkylsilyl, or CN; R11-R13 = independently H, halo, or alkyl; R14-R17 = independently H, halo, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)alkoxy, (halo)alkenyloxy, (halo)alkynyloxy, (halo)alkylthio, cycloalkyl(oxy), trialkylsilyl, or CN; X1 = O or S; X2 = O or S] are prepared as agrochem. fungicides. Thus, N-(3',4'-dimethoxybiphenyl-2-yl)-2-hydroxy-2-(4-methylphenyl)acetamide (preparation given) was treated with MeSO₂Cl in Et₂O in the presence of Et₃N, followed by the addition of 2-propyn-1-ol to give II. I showed antifungal activity of >90% at the concentration of 500 ppm in 4 days in tomato seedlings.

IT 602280-68-2P 602280-69-3P 602280-70-6P
602280-71-7P 602280-72-8P 602280-73-9P
602280-74-0P 602280-75-1P 602280-76-2P
602280-77-3P 602280-79-5P 602280-84-2P

(fungicide; preparation of N-biphenylacetamides as agrochem. fungicides)

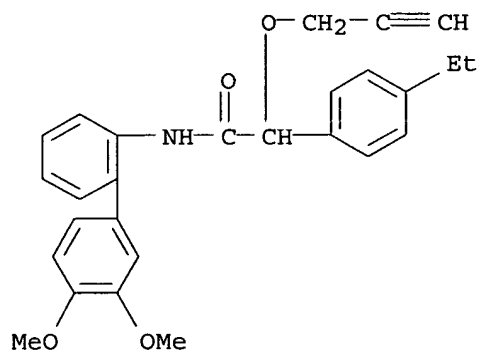
RN 602280-68-2 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-4-methyl- α -(2-propynyloxy)- (9CI) (CA INDEX NAME)



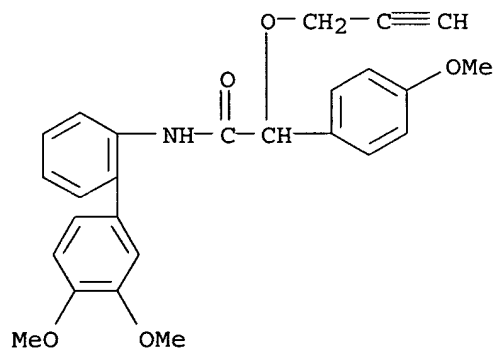
RN 602280-69-3 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-4-ethyl-
α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



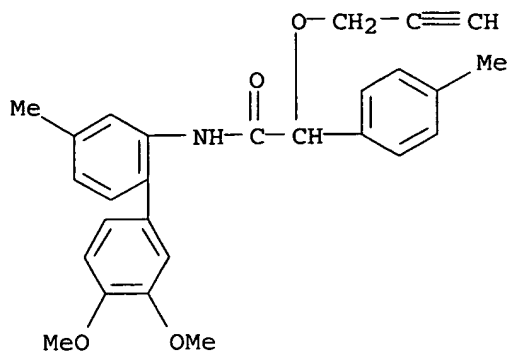
RN 602280-70-6 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-4-
methoxy-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)

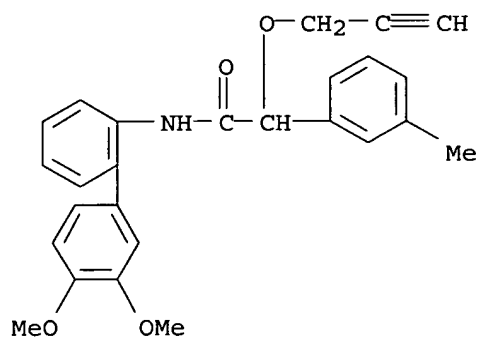


RN 602280-71-7 HCAPLUS

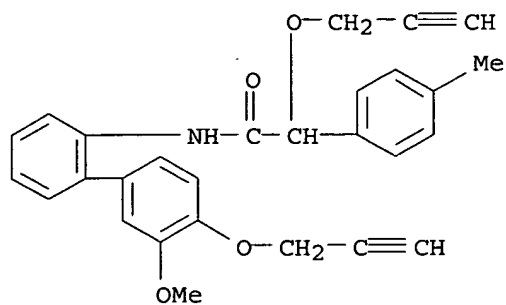
CN Benzeneacetamide, N-(3',4'-dimethoxy-4-methyl[1,1'-biphenyl]-2-yl)-
4-methyl-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



RN 602280-72-8 HCAPLUS

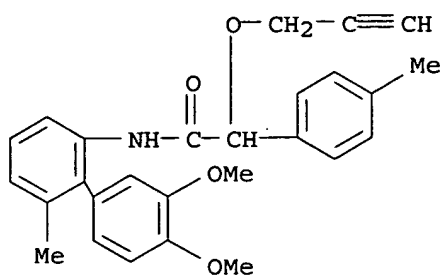
CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-3-methyl-
α-(2-propynyloxy)- (9CI) (CA INDEX NAME)

RN 602280-73-9 HCAPLUS

CN Benzeneacetamide, N-[3'-methoxy-4'-(2-propynyloxy)[1,1'-biphenyl]-
2-yl]-4-methyl-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)

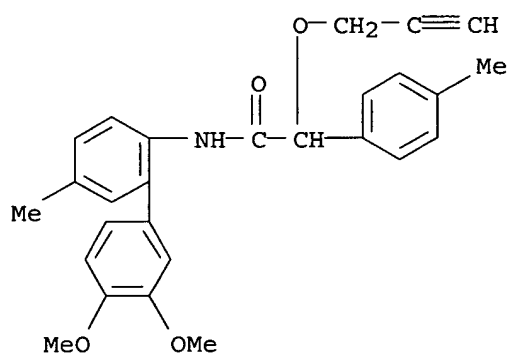
RN 602280-74-0 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy-6-methyl[1,1'-biphenyl]-2-yl)-
4-methyl-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



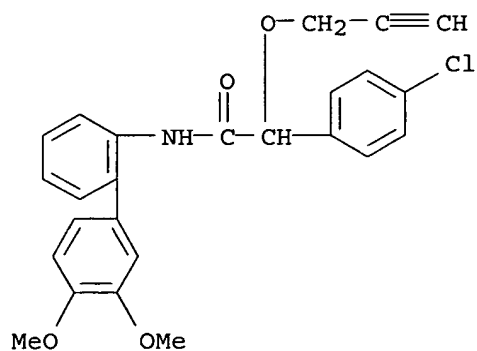
RN 602280-75-1 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy-5-methyl[1,1'-biphenyl]-2-yl)-4-methyl-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



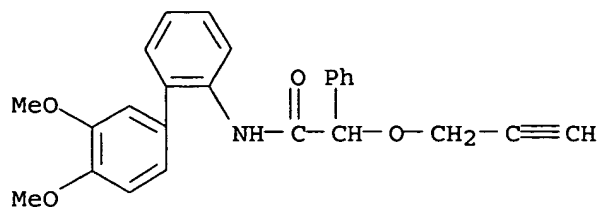
RN 602280-76-2 HCAPLUS

CN Benzeneacetamide, 4-chloro-N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



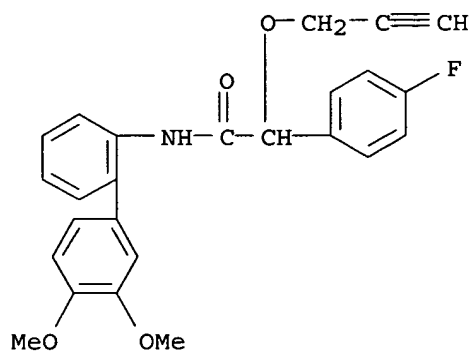
RN 602280-77-3 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



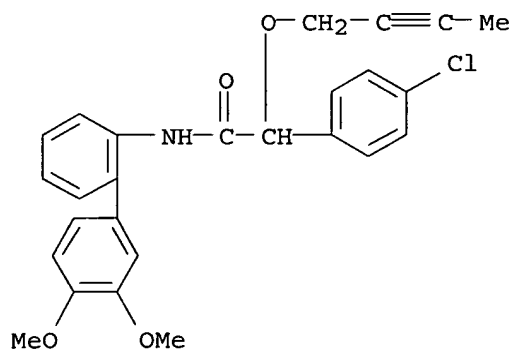
RN 602280-79-5 HCAPLUS

CN Benzeneacetamide, N-(3',4'-dimethoxy[1,1'-biphenyl]-2-yl)-4-fluoro-
 α -(2-propynyloxy)- (9CI) (CA INDEX NAME)



RN 602280-84-2 HCAPLUS

CN Benzeneacetamide, α -(2-butynyloxy)-4-chloro-N-(3',4'-
dimethoxy[1,1'-biphenyl]-2-yl)- (9CI) (CA INDEX NAME)



IC ICM C07C235-38

ICS C07C255-13; C07C255-63; C07C323-37; C07C323-60; C07C323-62;
C07C327-44

CC 25-19 (Benzene, Its Derivatives, and Condensed Benzenoid
Compounds)

Section cross-reference(s): 5

IT 602280-68-2P 602280-69-3P 602280-70-6P

602280-71-7P 602280-72-8P 602280-73-9P

602280-74-0P 602280-75-1P 602280-76-2P

602280-77-3P 602280-78-4P 602280-79-5P

602280-80-8P 602280-81-9P 602280-82-0P 602280-84-2P

602280-86-4P 602280-87-5P 602280-88-6P 602280-89-7P

602281-06-1P

(fungicide; preparation of N-biphenylacetamides as agrochem.
fungicides)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE
FOR THIS RECORD. ALL CITATIONS AVAILABLE
IN THE RE FORMAT

L16 ANSWER 3 OF 3 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2001:900234 HCAPLUS

DOCUMENT NUMBER: 136:340462

TITLE: Synthesis and biological activity of novel
thyroid hormone analogues: 5'-aryl substituted
GC-1 derivatives

AUTHOR(S): Chiellini, Grazia; Nguyen, Ngoc-Ha; Apriletti,
James W.; Baxter, John D.; Scanlan, Thomas S.

CORPORATE SOURCE: Departments of Pharmaceutical Chemistry and
Cellular & Molecular Pharmacology, University
of California, San Francisco, CA, 94143-0446,
USA

SOURCE: Bioorganic & Medicinal Chemistry (2001),
Volume Date 2002, 10(2), 333-346
CODEN: BMECEP; ISSN: 0968-0896

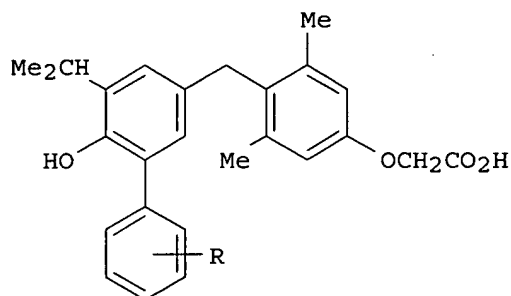
PUBLISHER: Elsevier Science Ltd.

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 136:340462

GI



I

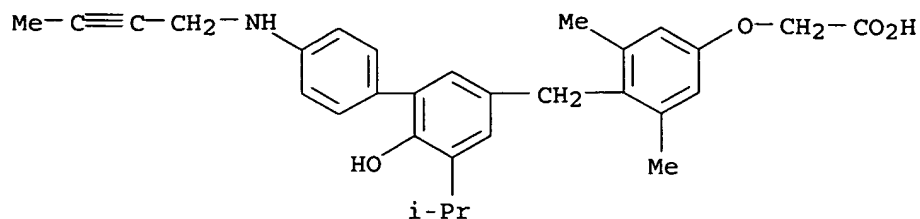
AB Biphenylmethylphenoxyacetic acids I [R = 4-NO₂, 4-NHCH₂CO₂H,
4-NHCONHPh, 4-NHCH₂C.tplbond.CMe, 4-NH₂, 3-NO₂, 2-NO₂, 4-CO₂H,
4-CONH₂, 4-NHC(:NH)NH₂] were prepared via arylation of the
diphenylmethaneboronic acid. Substitution at the 5'-position
decreased binding affinity, but retained TR β -selectivity for
most of the compds. Transactivation assays reveal that most of
these compds. function as thyroid hormone agonists, but I [R =
4-NO₂] antagonizes the response to thyroid hormone.

IT 417872-45-8P

(preparation of biphenylmethylphenoxyacetic acids as thyroid hormone
analogues)

RN 417872-45-8 HCAPLUS

CN Acetic acid, [4-[[4'-(2-butynylamino)-6-hydroxy-5-(1-
methylethyl)[1,1'-biphenyl]-3-yl]methyl]-3,5-dimethylphenoxy]-
(9CI) (CA INDEX NAME)



CC 25-17 (Benzene, Its Derivatives, and Condensed Benzenoid Compounds)

Section cross-reference(s): 1

IT 417871-97-7P 417872-05-0P 417872-10-7P 417872-14-1P
 417872-18-5P 417872-30-1P 417872-38-9P 417872-45-8P
 417872-54-9P 417872-67-4P 447415-34-1P

(preparation of biphenylmethylphenoxyacetic acids as thyroid hormone analogs)

REFERENCE COUNT: 30 THERE ARE 30 CITED REFERENCES AVAILABLE
 FOR THIS RECORD. ALL CITATIONS AVAILABLE
 IN THE RE FORMAT

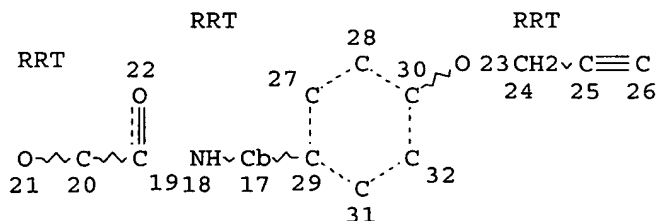
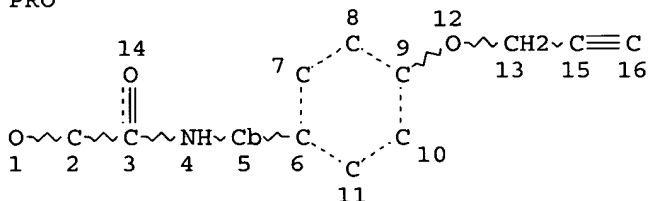
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=> d que stat l19

L17 STR

PRO



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GGCAT IS MCY UNS AT 5

GGCAT IS MCY UNS AT 17

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 32

STEREO ATTRIBUTES: NONE

L19 0 SEA FILE=CASREACT SSS FUL L17 (0 REACTIONS)

100.0% DONE 257 VERIFIED 0 HIT RXNS 0 DO

CS

SEARCH TIME: 00.00.01